C. Amendments to the Claims

Please amend the claims where indicated:

1. (original) A compound having the formula

(I)

wherein

- R₁ and R₂ are independently hydrogen; lower alkyl; C₁-C₆ cycloalkyl or cycloheteroalkyl; halogen or halo-substituted alkyl; or R₁ and R₂, taken together, form a C₅-C₇ cycloalkyl or cycloheteroalkyl ring;
- Cy is a single or conjugated substituted or unsubstituted alicyclic or aromatic ring structure; and
- n is 0, 1, 2, 3, 4 or 5; and pharmaceutically acceptable salts and/or esters thereof.
- 2. (original) The compound of claim 1, wherein R₁ and R₂, taken together, form a C₅-C₇ cycloalkyl or cycloheteroalkyl ring.
- 3. (original) The compound of claim 2, wherein R₁ and R₂, taken together, form a cyclohexyl ring.
- 4. (original) The compound of claim 1, wherein n is 0, 1, 2 or 3.
- 5. (original) The compound of claim 1, wherein said lower alkyl is C_1 - C_5 alkyl.
- (original) The compound of claim 1, wherein said compound is a 5-HT receptor antagonist.
- 7. (original) The compound of claim 6, wherein said compound is a 5-HT₂ receptor antagonist.
- 8. (original) The compound of claim 7, wherein said compound is a 5-HT_{2A, B or C} receptor antagonist.

- 9. (original) The compound of claim 7, wherein said compound is a 5-HT_{2B} receptor antagonist.
- 10. (original) The compound of claim 1, wherein said compound is [1-(2-Fluoro-benzyl)-piperidin-4-yl]-(5,6,7,8-tetrahydro-benzo[4,5]thieno[2,3-d]pyrimidin-4-yl)-amine.
- 11. (original) The compound of claim 1, wherein said compound is [1-(3-Fluoro-benzyl)-piperidin-4-yl]-(5,6,7,8-tetrahydro-benzo[4,5]thieno[2,3-d]pyrimidin-4-yl)-amine.
- 12. (original) The compound of claim 1, wherein said compound is [1-(4-Fluoro-benzyl)-piperidin-4-yl]-(5,6,7,8-tetrahydro-benzo[4,5]thieno[2,3-d]pyrimidin-4-yl)-amine.
- 13. (original) The compound of claim 1, wherein said compound is [1-(4-Methyl-benzyl)-piperidin-4-yl]-(5,6,7,8-tetrahydro-benzo[4,5]thieno[2,3-d]pyrimidin-4-yl)-amine.
- 14. (original) The compound of claim 1, wherein said compound is (1-Benzo[1,3]dioxol-5-ylmethyl-piperidin-4-yl)-(5,6,7,8-tetrahydro-benzo[4,5]thieno[2,3-d]pyrimidin-4-yl)-amine.
- 15. (original) The compound of claim 1, wherein said compound is (5,6,7,8-Tetrahydrobenzo[4,5]thieno[2,3-d]pyrimidin-4-yl)-[1-(4-trifluoromethyl-benzyl)-piperidin-4-yl]-amine.
- 16. (original) The compound of claim 1, wherein said compound is (1-Benzhydryl-piperidin-4-yl)-(5,6,7,8-tetrahydro-benzo[4,5]thieno[2,3-d]pyrimidin-4-yl)-amine.
- 17. (original) The compound of claim 1, wherein said compound is (1-Naphthalen-2-ylmethyl-piperidin-4-yl)-(5,6,7,8-tetrahydro-benzo[4,5]thieno[2,3-d]pyrimidin-4-yl)-amine.
- 18. (original) The compound of claim 1, wherein said compound is (1-Phenethyl-piperidin-4-yl)-(5,6,7,8-tetrahydro-benzo[4,5]thieno[2,3-d]pyrimidin-4-yl)-amine.
- 19. (original) The compound of claim 1, wherein said compound is [1-(3-Phenyl-propyl)-piperidin-4-yl]-(5,6,7,8-tetrahydro-benzo[4,5]thieno[2,3-d]pyrimidin-4-yl)-amine.
- 20. (original) The compound of claim 1, wherein said compound is (5,6,7,8-Tetrahydrobenzo[4,5]thieno[2,3-d]pyrimidin-4-yl)-[1-(4-trifluoromethoxy-benzyl)-piperidin-4-yl]-amine.

- 21. (original) The compound of claim 1, wherein said compound is [1-(4-Methoxy-benzyl)-piperidin-4-yl]-(5,6,7,8-tetrahydro-benzo[4,5]thieno[2,3-d]pyrimidin-4-yl)-amine.
- 22. (original) A compound having the formula

$$\begin{array}{c|c} & & & & \\ & & & & \\ & & & & \\ R_1 & & & \\ & & & \\ R_2 & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & \\ & & & \\ &$$

wherein

- R₁ and R₂ are independently be hydrogen; lower alkyl, C₁-C₆ cycloalkyl or cycloheteroalkyl; halogens or halo-substituted alkyl; or R₁ and R₂, taken together, form a C₅-C₇ cycloalkyl or cycloheteroalkyl ring;
- R₃ and R₄ are independently Ar which is a single or conjugated substituted or unsubstituted aromatic ring structure;
- R₅ is H, (C₁-C₅)alkyl, (C₁-C₆)cycloalkyl, halogen substituted alkyl, NH₂,
 NHMe, NMe₂, NHEt, NH(Et)₂, NH(Pr), N(Pr)₂; and
- n is 0, 1, 2, 3, 4 or 5; and pharmaceutically acceptable salts and/or esters thereof.

23-32. (canceled)